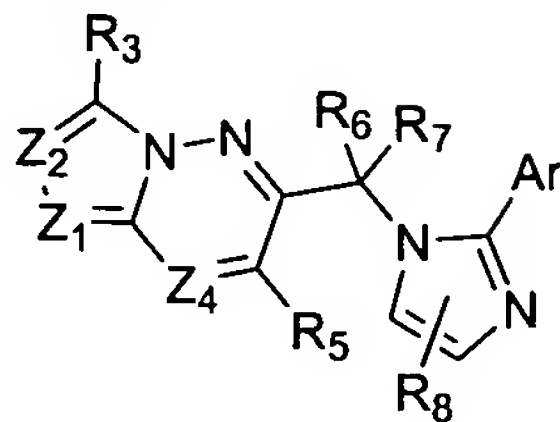


AMENDMENTS TO THE CLAIMS

1. (Original) A compound of the formula:



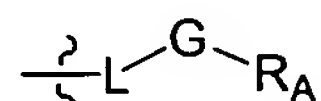
or a pharmaceutically acceptable salt thereof, wherein:

Z₁ is nitrogen or CR₁ and Z₂ is nitrogen or CR₂; such that at least one of Z₁ and Z₂ is nitrogen;

Z₄ is nitrogen or CR₄;

R₁, R₂, R₃ and R₄ are each independently selected from:

- (a) hydrogen, halogen, nitro and cyano; and
- (b) groups of the formula:



wherein:

L is a single covalent bond or C₁-C₈alkylene;

G is a single covalent bond, N(R_B), O, C(=O), C(=O)O, C(=O)N(R_B), N(R_B)C(=O), S(O)_m, CH₂C(=O), S(O)_mN(R_B) or N(R_B)S(O)_m; wherein m is 0, 1 or 2; and

R_A and each R_B are independently selected from:

- (i) hydrogen; and
- (ii) C₁-C₈alkyl, C₂-C₈alkenyl, C₂-C₈alkynyl, (C₃-C₈cycloalkyl)C₀-C₄alkyl, (3- to 6-membered heterocycloalkyl)C₀-C₄alkyl, (aryl)C₀-C₂alkyl and (heteroaryl)C₀-C₂alkyl, each of which is substituted with from 0 to 4 substituents independently selected from halogen, hydroxy, nitro, cyano, amino, C₁-C₄alkyl, C₁-C₄alkoxy, C₁-C₄alkanoyl, mono- and di(C₁-C₄alkyl)amino, C₁-C₄haloalkyl and C₁-C₄haloalkoxy;

R₅ is hydrogen, halogen, cyano, C₁-C₆alkyl, C₂-C₆alkenyl, C₂-C₆alkynyl, C₁-C₄alkoxy, or mono- or di-(C₁-C₄alkyl)amino, each of which is substituted with from 0 to 5 substituents independently chosen from halogen, hydroxy, nitro, cyano, amino, C₁-C₄alkoxy, C₁-C₂haloalkyl, C₁-C₂haloalkoxy, mono- and di-C₁-C₄alkylamino, C₃-C₈cycloalkyl, phenyl, phenylC₁-C₄alkoxy and 5- or 6-membered heteroaryl;

R₆ and R₇ are independently hydrogen, methyl, ethyl or halogen;

R₈ represents 0, 1 or 2 substituents independently chosen from halogen, hydroxy, nitro, cyano, amino, C₁-C₄alkyl, C₁-C₄alkoxy, mono- and di(C₁-C₄alkyl)amino, C₃-C₇cycloalkyl, C₁-C₂haloalkyl and C₁-C₂haloalkoxy; and

Ar represents phenyl, naphthyl or a 5- to 10-membered heteroaryl, each of which is substituted with from 0 to 4 substituents independently chosen from halogen, hydroxy, nitro, cyano, amino, C₁-

C₈alkyl, C₁-C₈alkenyl, C₁-C₈alkynyl, C₁-C₈alkoxy, (C₃-C₇cycloalkyl)C₀-C₄alkyl, (C₃-C₇cycloalkyl)C₁-C₄alkoxy, C₂-C₈alkyl ether, C₃-C₈alkanone, C₁-C₈alkanoyl, 3- to 7-membered heterocycloalkyl, C₁-C₈haloalkyl, C₁-C₈haloalkoxy, oxo, C₁-C₈hydroxyalkyl, C₁-C₈aminoalkyl, and mono- and di-(C₁-C₈alkyl)aminoC₀-C₈alkyl.

2. (Original) A compound or salt according to claim 1, wherein R₈ represents 0 or 1 substituent selected from halogen, C₁-C₂alkyl and C₁-C₂alkoxy.

3. (Currently Amended) A compound or salt according to claim 1 ~~or claim 2~~, wherein Ar is substituted with 0, 1, 2 or 3 substituents independently selected from halogen, hydroxy, amino, cyano, C₁-C₄alkyl, C₁-C₄alkoxy, mono- or di-C₁-C₄alkylamino, C₂-C₄alkanoyl, (C₃-C₇cycloalkyl)C₀-C₂alkyl, C₁-C₂haloalkyl and C₁-C₂haloalkoxy.

4. (Currently Amended) A compound or salt according to claim 1 ~~or claim 2~~, wherein Ar represents phenyl, pyridyl, thiazolyl, thienyl, pyridazinyl or pyrimidinyl, each of which is substituted with from 0 to 4 substituents.

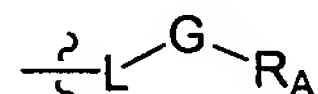
5. (Original) A compound or salt according to claim 4, wherein Ar represents phenyl, pyridyl, thiazolyl, thienyl or pyridazinyl, each of which is substituted with from 0 to 3 substituents independently selected from chloro, fluoro, hydroxy, cyano, amino, C₁-C₄alkyl, C₁-C₄alkoxy, C₁-C₂alkylamino, C₁-C₂haloalkyl and C₁-C₂haloalkoxy.

6. (Original) A compound or salt according to claim 5, wherein Ar represents phenyl, 2-pyridyl, 1,3-thiazol-2-yl, 2-thienyl or pyridazin-3-yl, each of which is substituted with from 0 to 3 substituents independently selected from fluoro, chloro, hydroxy, C₁-C₂alkyl, cyano and C₁-C₂alkoxy.

7. (Original) A compound or salt according to claim 5, wherein Ar represents 2,6-difluoro-phenyl, 2,5-difluoro-phenyl, 5-fluoro-2-methyl-phenyl, pyridine-2-yl, 3-fluoro-pyridin-2-yl, 3-cyano-pyridin-2-yl, 3-trifluoromethyl-pyridin-2-yl, 3-hydroxy-pyridin-2-yl, 3-methoxy-pyridin-2-yl, 6-fluoro-pyridin-2-yl, 6-cyano-pyridin-2-yl, 6-trifluoromethyl-pyridin-2-yl, 6-hydroxy-pyridin-2-yl or 6-methoxy-pyridin-2-yl.

8. (Currently Amended) A compound or salt according to ~~any one of claims 1-7~~ claim 1, wherein R₁, R₂, R₃ and R₄ are independently selected from:

- (a) hydrogen, halogen or cyano; and
- (b) groups of the formula:



wherein:

- (i) L is a single covalent bond;
- (ii) G is a single covalent bond, -NH-, -N(R_B)-, -O-, -C(=O)O- or C(=O)-; and
- (iii) R_A and R_B are independently selected from (1) hydrogen and (2) C₁-C₆alkyl, C₂-C₆alkenyl, (C₃-C₇cycloalkyl)C₀-C₂alkyl, phenyl, thienyl, pyridyl, pyrimidinyl, thiazolyl and pyrazinyl, each of which is substituted with from 0 to 4 substituents independently selected from hydroxy, halogen, cyano, amino, C₁-C₂alkyl and C₁-C₂alkoxy.

9. (Original) A compound or salt according to claim 8 wherein R₁, R₂, R₃ and R₄ are independently selected from hydrogen, hydroxy, halogen, cyano, aminocarbonyl, C₁-C₆alkyl, C₁-C₆alkoxy, C₂-C₆alkyl ether, C₃-C₇cycloalkyl, C₁-C₄hydroxyalkyl, C₁-C₂haloalkyl, C₁-C₂haloalkoxy, C₁-C₆alkoxycarbonyl, mono- and di-(C₁-C₄alkyl)amino, phenyl and pyridyl.

10. (Original) A compound or salt according to claim 9, wherein R₃ and R₄ are independently selected from hydrogen, methyl and ethyl.

11. (Currently Amended) A compound or salt according to ~~any one of claims 1-10,~~ claim 8 wherein Z₁ is nitrogen and Z₂ is CR₂.

12. (Original) A compound or salt according to claim 11, wherein R₂ is selected from hydrogen, cyano, aminocarbonyl, C₁-C₄alkyl, C₁-C₄alkoxy, C₁-C₄alkoxycarbonyl, C₂-C₄alkyl ether, C₃-C₇cycloalkyl, C₁-C₂hydroxyalkyl, fluoromethyl, difluoromethyl, trifluoromethyl, phenyl and pyridyl.

13. (Currently Amended) A compound or salt according to ~~any one of claims 1-10,~~ claim 8, wherein Z₁ is CR₁ and Z₂ is nitrogen.

14. (Original) A compound or salt according to claim 13, wherein R₁ is selected from hydrogen, cyano, aminocarbonyl, C₁-C₄alkyl, C₁-C₄alkoxy, C₁-C₄alkoxycarbonyl, C₂-C₄alkyl ether, C₃-C₇cycloalkyl, C₁-C₂hydroxyalkyl, fluoromethyl, difluoromethyl, trifluoromethyl, phenyl and pyridyl.

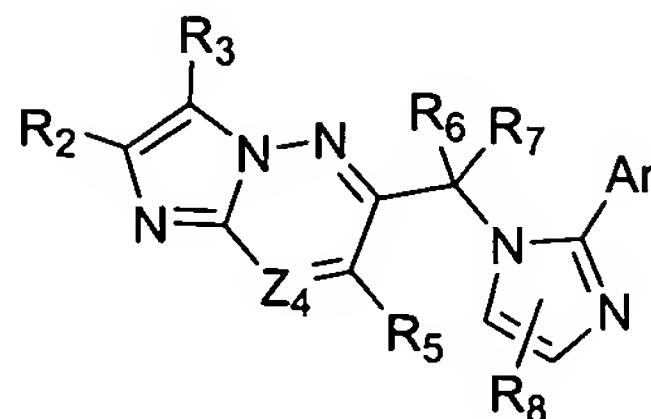
15. (Currently Amended) A compound or salt according to ~~any one of claims 1-10,~~
claim 8, wherein Z₁ and Z₂ are nitrogen.

16. (Currently Amended) A compound or salt according to ~~any one of claims 1-15,~~
claim 8, wherein R₆ and R₇ are both hydrogen.

17. (Currently Amended) A compound or salt according to ~~any one of claims 1-16,~~
claim 8 wherein R₅ is C₁-C₆alkyl, C₂-C₆ alkenyl, C₁-C₄ alkoxy, or mono- or di-C₁-C₄alkylamino, each of which is substituted with from 0 to 2 substituents independently selected from halogen, hydroxy, C₁-C₂alkoxy, C₃-C₈cycloalkyl, phenyl and phenylC₁-C₂alkoxy.

18. (Original) A compound or salt according to claim 17, wherein R₅ is ethyl, propyl, butyl, ethoxy or methoxymethyl.

19. (Original) A compound or salt according to claim 1, wherein the compound has the formula:



wherein:

R₂ is selected from hydrogen, hydroxy, halogen, cyano, aminocarbonyl, C₁-C₆alkyl, C₁-C₆alkoxy, C₃-C₇cycloalkyl, C₂-C₆alkyl ether, C₁-C₄hydroxyalkyl, C₁-C₂haloalkyl, C₁-C₂haloalkoxy, C₁-C₄alkoxycarbonyl, mono- and di-(C₁-C₄alkyl)amino, phenyl and pyridyl;

R₃ and R₄, if present, are independently hydrogen or C₁-C₄alkyl;

R₅ is C₁-C₆alkyl, C₂-C₆alkenyl, C₁-C₄alkoxy, or mono- or di-C₁-C₄alkylamino, each of which is substituted with from 0 to 2 substituents independently selected from halogen, hydroxy, C₁-C₂alkoxy, C₃-C₈cycloalkyl, phenyl and phenylC₁-C₂alkoxy;

R₆ and R₇ are independently hydrogen, methyl, ethyl or halogen;

R₈ represents 0 or 1 substituent selected from halogen, C₁-C₂alkyl and C₁-C₂alkoxy; and

Ar represents phenyl, 2-pyridyl, 1,3-thiazol-2-yl, 2-thienyl or 3-pyridazinyl, each of which is substituted with from 0 to 3 substituents independently selected from fluoro, chloro, hydroxy, C₁-C₂alkyl, C₁-C₂haloalkyl, cyano and C₁-C₂alkoxy.

20. (Original) A compound or salt according to claim 19, wherein:

R₂ is hydrogen, cyano, aminocarbonyl, C₁-C₆alkyl, C₁-C₆alkoxy, C₃-C₆cycloalkyl, C₂-C₆alkyl ether, C₁-C₄hydroxyalkyl, C₁-C₂haloalkyl or C₁-C₄alkoxycarbonyl;

Z₄ is CR₄;

R₃ and R₄ are independently hydrogen or C₁-C₂alkyl;

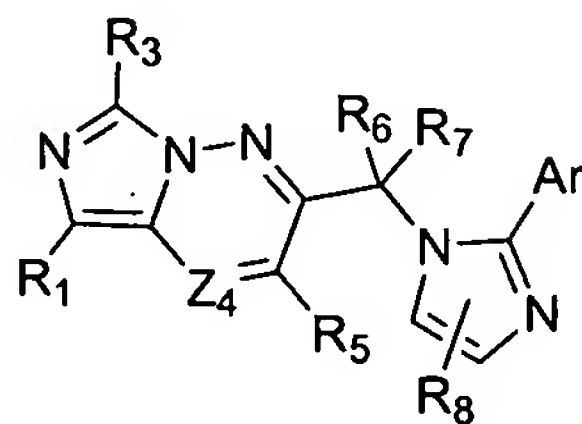
R₅ is C₁-C₆alkyl or C₂-C₆alkenyl, each of which is substituted with from 0 to 2 substituents independently selected from halogen, hydroxy and C₁-C₂alkoxy;

R₆ and R₇ are hydrogen;

R₈ represents 0 substituents; and

Ar represents phenyl or 2-pyridyl, each of which is substituted with from 0 to 3 substituents independently selected from fluoro, chloro, C₁-C₂alkyl, C₁-C₂haloalkyl, cyano and C₁-C₂alkoxy.

21. (Original) A compound or salt according to claim 1, wherein the compound has the formula:



wherein:

R₁ is selected from hydrogen, hydroxy, halogen, cyano, aminocarbonyl, C₁-C₆alkyl, C₁-C₆alkoxy, C₃-C₇cycloalkyl, C₂-C₆alkyl ether, C₁-C₄hydroxyalkyl, C₁-C₂haloalkyl, C₁-C₂haloalkoxy, C₁-C₄alkoxycarbonyl, mono- and di-(C₁-C₄alkyl)amino, phenyl and pyridyl;

R₃ and R₄, if present, are independently hydrogen or C₁-C₄alkyl;

R₅ is C₁-C₆ alkyl, C₂-C₆ alkenyl, C₁-C₄ alkoxy, or mono- or di-C₁-C₄alkylamino, each of which is substituted with from 0 to 2 substituents independently selected from halogen, hydroxy, C₁-C₂alkoxy, C₃-C₈cycloalkyl, phenyl and phenylC₁-C₂alkoxy;

R₆ and R₇ are independently hydrogen, methyl, ethyl or halogen;

R₈ represents 0 or 1 substituent selected from halogen, C₁-C₂alkyl and C₁-C₂alkoxy; and

Ar represents phenyl, 2-pyridyl, 1,3-thiazol-2-yl, 2-thienyl or 3-pyridazinyl, each of which is substituted with from 0 to 3 substituents independently selected from fluoro, chloro, hydroxy, C₁-C₂alkyl, C₁-C₂haloalkyl, cyano and C₁-C₂alkoxy.

22. (Original) A compound or salt according to claim 21, wherein:

R₁ is hydrogen, cyano, aminocarbonyl, C₁-C₆alkyl, C₁-C₆alkoxy, C₃-C₆cycloalkyl, C₂-C₆alkyl ether, C₁-C₄hydroxyalkyl, C₁-C₂haloalkyl or C₁-C₄alkoxycarbonyl;

Z₄ is CR₄;

R₃ and R₄ are independently hydrogen or C₁-C₂alkyl;

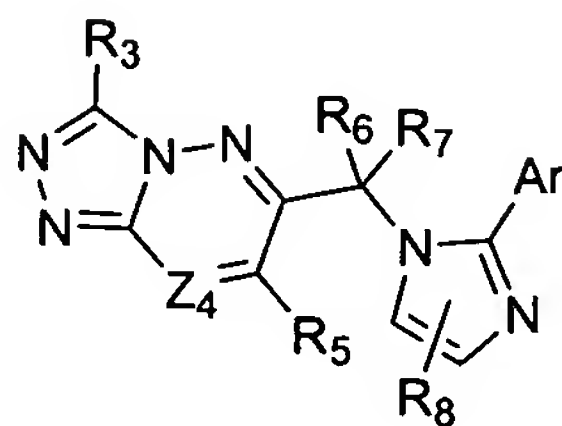
R₅ is C₁-C₆alkyl or C₂-C₆alkenyl, each of which is substituted with from 0 to 2 substituents independently selected from halogen, hydroxy and C₁-C₂alkoxy;

R₆ and R₇ are hydrogen;

R₈ represents 0 substituents; and

Ar represents phenyl or 2-pyridyl, each of which is substituted with from 0 to 3 substituents independently selected from fluoro, chloro, C₁-C₂alkyl, C₁-C₂haloalkyl, cyano and C₁-C₂alkoxy.

23. (Original) A compound or salt according to claim 1, wherein the compound has the formula:



wherein:

R₃ and R₄, if present, are independently hydrogen or C₁-C₄alkyl;

R₅ is C₁-C₆ alkyl, C₂-C₆ alkenyl, C₁-C₄ alkoxy, or mono- or di-C₁-C₄alkylamino, each of which is substituted with from 0 to 2 substituents independently selected from halogen, hydroxy, C₁-C₂alkoxy, C₃-C₈cycloalkyl, phenyl and phenylC₁-C₂alkoxy;

R₆ and R₇ are independently hydrogen, methyl, ethyl or halogen;

R₈ represents 0 or 1 substituent selected from halogen, C₁-C₂alkyl and C₁-C₂alkoxy; and

Ar represents phenyl, 2-pyridyl, 1,3-thiazol-2-yl, 2-thienyl or 3-pyridazinyl, each of which is substituted with from 0 to 3 substituents independently selected from fluoro, chloro, hydroxy, C₁-C₂alkyl, C₁-C₂haloalkyl, cyano and C₁-C₂alkoxy.

24. (Original) A compound or salt according to claim 23, wherein:

Z₄ is CR₄;

R₃ and R₄ are independently hydrogen or C₁-C₂alkyl;

R₅ is C₁-C₆alkyl or C₂-C₆alkenyl, each of which is substituted with from 0 to 2 substituents independently selected from halogen, hydroxy and C₁-C₂alkoxy;

R₆ and R₇ are hydrogen;

R₈ represents 0 substituents; and

Ar represents phenyl or 2-pyridyl, each of which is substituted with from 0 to 3 substituents independently selected from fluoro, chloro, C₁-C₂alkyl, C₁-C₂haloalkyl, cyano and C₁-C₂alkoxy.

25. (Currently Amended) A compound or salt according to ~~any one of claims 1-24~~ claim 1, wherein the compound exhibits a K_i of 1 micromolar or less in an assay of GABA_A receptor binding.

26. (Original) A compound or salt according to claim 25, wherein the compound exhibits a K_i of 100 nanomolar or less in an assay of GABA_A receptor binding.

27. (Original) A compound or salt according to claim 26, wherein the compound exhibits a K_i of 10 nanomolar or less in an assay of GABA_A receptor binding.

28. (Currently Amended) A pharmaceutical composition comprising a compound or salt according to ~~any one of claims 1-24~~ claim 1 in combination with a physiologically acceptable carrier or excipient.

29. (Original) A pharmaceutical composition according to claim 28, wherein the pharmaceutical composition is formulated as an injectible fluid, an aerosol, a cream, a gel, a pill, a capsule, a syrup or a transdermal patch.

30. (Currently Amended) A method for the treatment of anxiety, depression, or a sleep disorder, ~~attention-deficit disorder or Alzheimer's dementia~~, comprising administering to a patient in need of such treatment a therapeutically effective amount of a compound or salt according to ~~any one of claims 1-24~~ claim 1.

31 – 38. Cancelled.

39. (Currently Amended) A packaged pharmaceutical preparation comprising a pharmaceutical composition according to claim 28 in a container and instructions for using the composition to treat a patient suffering from anxiety, depression, a sleep disorder, ~~attention-deficit disorder, Alzheimer's dementia~~ or short-term memory loss.

40. Cancelled.

41. (Currently Amended) A compound or salt of Claim 1, wherein the compound is
6-[2-(6-fluoro-pyridin-2-yl)-imidazol-1-ylmethyl]-7-propyl-imidazo[1,2-b]pyridazine;
2-tert-butyl-6-[2-(6-fluoro-pyridin-2-yl)-imidazol-1-ylmethyl]-7-propyl-imidazo[1,2-b]pyridazine
2-ethyl-6-[2-(6-fluoro-pyridin-2-yl)-imidazol-1-ylmethyl]-7-propyl-imidazo[1,2-b]pyridazine
2-methyl-6-[2-(6-fluoro-pyridin-2-yl)-imidazol-1-ylmethyl]-7-propyl-imidazo[1,2-b]pyridazine
6-[2-(6-fluoro-pyridin-2-yl)-imidazol-1-ylmethyl]-7-propyl-2trifluoromethyl-imidazo[1,2-
b]pyridazine
6-[2-(6-fluoro-pyridin-2-yl)-imidazol-1-ylmethyl]-7-propyl-imidazo[1,2-b]pyridazine-2-carboxylic
acid ethyl ester
6-[2-(6-fluoro-pyridin-2-yl)-imidazol-1-ylmethyl]-7-propyl-imidazo[1,2-b]pyridazine-2-carboxylic
acid amide;
6-[2-(6-fluoro-pyridin-2-yl)-imidazol-1-ylmethyl]-7-propyl-imidazo[1,2-b]pyridazine-2-carbonitrile;
6-[2-(6-Fluoro-pyridin-2-yl)-imidazol-1-ylmethyl]-8-methyl-7-propyl-imidazo[1,2-b]pyridazine;
2-[2-(6-fluoro-pyridin-2-yl)-imidazol-1-ylmethyl]-5-methyl-3-propyl-imidazo[1,5-b]pyridazine;
7-ethyl-6-[2-(3-fluoro-phenyl)-imidazol-1-ylmethyl]-[1,2,4]triazolo[4,3-b] pyridazine; or
7-ethyl-6-[2-(3-fluoro-phenyl)-imidazol-1-ylmethyl]-3-methyl-[1,2,4]triazolo [4,3-b] pyridazine or a
pharmaceutically acceptable salt thereof.

42 – 52. Cancelled.